

Method Path : Z:\svoasrv\HPCHEM1\BNA_F\Methods\
 Method File : 8270-BF091422.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 Last Update : Wed Sep 14 14:55:51 2022
 Response Via : Initial Calibration

Calibration Files

5 =BF130232.D 10 =BF130233.D 20 =BF130234.D 40 =BF130235.D 50 =BF130236.D 60 =BF130237.D 80 =BF130238.D

	Compound	5	10	20	40	50	60	80	Avg	%RSD
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1) I	1,4-Dichlorobenzene	-----	-----	-----	-----	-----	-----	-----	-----	-----
2)	1,4-Dioxane	0.575	0.535	0.561	0.529	0.509	0.517	0.481	0.530	6.00
3)	Pyridine	1.410	1.410	1.509	1.338	1.342	1.366	1.296	1.381	5.01
4)	n-Nitrosodimethylamine	0.746	0.742	0.800	0.758	0.742	0.757	0.714	0.751	3.46
5) S	2-Fluorophenol	1.191	1.171	1.244	1.136	1.117	1.139	1.073	1.153	4.79
6)	Aniline	1.871	1.840	1.905	1.753	1.715	1.727	1.633	1.778	5.48
7) S	Phenol-d6	1.528	1.478	1.540	1.422	1.397	1.407	1.326	1.443	5.33
8)	2-Chlorophenol	1.350	1.343	1.425	1.295	1.275	1.287	1.220	1.314	5.00
9)	Benzaldehyde	1.134	1.078	1.106	0.933	0.895	0.855	0.765	0.966	14.59
10) C	Phenol	1.766	1.729	1.846	1.682	1.632	1.640	1.541	1.691	5.90
11)	bis(2-Chloroethyl)ether	1.302	1.266	1.319	1.206	1.169	1.169	1.106	1.220	6.44
12)	1,3-Dichlorobenzene	1.532	1.464	1.544	1.414	1.408	1.419	1.336	1.445	5.09
13) C	1,4-Dichlorobenzene	1.572	1.489	1.578	1.445	1.432	1.443	1.358	1.474	5.37
14)	1,2-Dichlorobenzene	1.494	1.404	1.485	1.362	1.317	1.315	1.261	1.377	6.45
15)	Benzyl Alcohol	1.091	1.103	1.266	1.174	1.144	1.148	1.081	1.144	5.58
16)	2,2'-oxybis(1,4-phenylene)	1.938	1.875	1.935	1.750	1.698	1.687	1.567	1.779	7.95
17)	2-Methylphenol	1.117	1.080	1.158	1.069	1.028	1.034	0.988	1.068	5.38
18)	Hexachloroethane	0.590	0.587	0.631	0.580	0.568	0.573	0.539	0.581	4.76
19) P	n-Nitroso-di-n-butylamine	1.033	0.999	1.050	0.978	0.933	0.928	0.895	0.974	5.91
20)	3+4-Methylphenols	1.490	1.429	1.505	1.382	1.316	1.331	1.256	1.387	6.69
21) I	Naphthalene-d8	-----	-----	-----	-----	-----	-----	-----	-----	-----
22)	Acetophenone	0.505	0.491	0.511	0.487	0.471	0.488	0.470	0.489	3.20
23) S	Nitrobenzene-d5	0.380	0.378	0.406	0.397	0.392	0.400	0.388	0.391	2.63
24)	Nitrobenzene	0.400	0.391	0.409	0.399	0.391	0.405	0.389	0.398	1.97
25)	Isophorone	0.620	0.628	0.650	0.633	0.621	0.643	0.622	0.631	1.85
26) C	2-Nitrophenol	0.138	0.149	0.167	0.171	0.168	0.177	0.171	0.163	8.73
27)	2,4-Dimethylphenol	0.238	0.250	0.258	0.253	0.250	0.258	0.249	0.251	2.73
28)	bis(2-Chloroethyl)ether	0.361	0.354	0.371	0.351	0.346	0.359	0.345	0.355	2.58
29) C	2,4-Dichlorophenol	0.262	0.266	0.281	0.279	0.273	0.287	0.277	0.275	3.09
30)	1,2,4-Trichlorobenzene	0.294	0.293	0.306	0.296	0.294	0.303	0.295	0.297	1.72
31)	Naphthalene	1.068	1.037	1.068	1.022	0.996	1.025	0.997	1.030	2.88
32)	Benzoic acid	0.141	0.179	0.209	0.206	0.217	0.213	0.194	0.194	15.08
33)	4-Chloroaniline	0.388	0.389	0.410	0.393	0.383	0.402	0.378	0.392	2.75
34) C	Hexachlorobutane	0.184	0.189	0.194	0.189	0.188	0.196	0.190	0.190	2.07
35)	Caprolactam	0.069	0.075	0.081	0.082	0.080	0.085	0.080	0.079	6.79
36) C	4-Chloro-3-methylphenol	0.293	0.295	0.320	0.309	0.304	0.312	0.300	0.305	3.14
37)	2-Methylnaphthalene	0.678	0.661	0.678	0.654	0.638	0.653	0.637	0.657	2.59
38)	1-Methylnaphthalene	0.641	0.642	0.657	0.625	0.610	0.631	0.611	0.631	2.70

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39) I	Acenaphthene-d10	-----ISTD-----	
40)	1,2,4,5-Tetrac...	0.544 0.526 0.553 0.531 0.539 0.564 0.564 0.546	2.73
41) P	Hexachlorocycl...	0.225 0.255 0.290 0.304 0.313 0.331 0.328 0.292	13.40
42) S	2,4,6-Tribromo...	0.170 0.182 0.188 0.195 0.193 0.208 0.205 0.192	6.77
43) C	2,4,6-Trichlor...	0.353 0.358 0.393 0.386 0.387 0.399 0.391 0.381	4.77
44)	2,4,5-Trichlor...	0.385 0.381 0.412 0.409 0.422 0.434 0.435 0.411	5.22
45) S	2-Fluorobiphenyl	1.385 1.366 1.428 1.346 1.355 1.386 1.349 1.373	2.11
46)	1,1'-Biphenyl	1.521 1.481 1.534 1.474 1.456 1.518 1.473 1.494	2.00
47)	2-Chloronaphth...	1.212 1.197 1.230 1.190 1.191 1.234 1.204 1.208	1.46
48)	2-Nitroaniline	0.353 0.371 0.415 0.417 0.416 0.429 0.419 0.403	7.14
49)	Acenaphthylene	1.854 1.791 1.871 1.784 1.782 1.838 1.761 1.812	2.32
50)	Dimethylphthalate	1.401 1.355 1.424 1.371 1.352 1.393 1.375 1.382	1.89
51)	2,6-Dinitrotol...	0.269 0.268 0.297 0.288 0.294 0.304 0.302 0.289	5.10
52) C	Acenaphthene	1.184 1.160 1.204 1.175 1.179 1.234 1.196 1.190	2.01
53)	3-Nitroaniline	0.299 0.298 0.325 0.328 0.326 0.342 0.335 0.322	5.26
54) P	2,4-Dinitrophenol	0.082 0.118 0.139 0.141 0.155 0.159 0.132	21.61
55)	Dibenzofuran	1.741 1.642 1.707 1.625 1.610 1.652 1.614 1.656	3.01
56) P	4-Nitrophenol	0.190 0.211 0.252 0.250 0.241 0.251 0.248 0.234	10.37
57)	2,4-Dinitrotol...	0.319 0.333 0.383 0.383 0.380 0.396 0.392 0.369	8.25
58)	Fluorene	1.363 1.337 1.376 1.355 1.326 1.372 1.349 1.354	1.35
59)	2,3,4,6-Tetrac...	0.314 0.293 0.334 0.319 0.322 0.338 0.333 0.322	4.77
60)	Diethylphthalate	1.370 1.355 1.425 1.389 1.375 1.404 1.381 1.385	1.67
61)	4-Chlorophenyl...	0.605 0.573 0.599 0.589 0.581 0.609 0.603 0.594	2.28
62)	4-Nitroaniline	0.294 0.303 0.339 0.331 0.330 0.336 0.331 0.323	5.47
63)	Azobenzene	1.446 1.424 1.485 1.430 1.391 1.424 1.386 1.427	2.35
64) I	Phenanthrene-d10	-----ISTD-----	
65)	4,6-Dinitro-2...	0.083 0.102 0.115 0.115 0.125 0.127 0.111	14.94
66) c	n-Nitrosodiphe...	0.671 0.658 0.691 0.655 0.654 0.678 0.671 0.668	2.00
67)	4-Bromophenyl...	0.212 0.210 0.224 0.216 0.224 0.230 0.229 0.221	3.68
68)	Hexachlorobenzene	0.243 0.239 0.257 0.245 0.247 0.259 0.259 0.250	3.27
69)	Atrazine	0.191 0.193 0.199 0.194 0.196 0.200 0.195 0.195	1.66
70) C	Pentachlorophenol	0.102 0.117 0.134 0.141 0.145 0.150 0.151 0.134	13.83
71)	Phenanthrene	1.163 1.136 1.157 1.095 1.090 1.128 1.092 1.123	2.76
72)	Anthracene	1.094 1.085 1.122 1.070 1.068 1.093 1.053 1.084	2.06
73)	Carbazole	1.010 0.972 1.017 0.970 0.956 0.962 0.958 0.978	2.59
74)	Di-n-butylphth...	1.117 1.155 1.256 1.192 1.178 1.185 1.172 1.179	3.58
75) C	Fluoranthene	1.116 1.098 1.162 1.086 1.046 1.051 1.036 1.085	4.13
76) I	Chrysene-d12	-----ISTD-----	
77)	Benzidine	0.629 0.570 0.593 0.615 0.560 0.601 0.496 0.581	7.61
78)	Pyrene	1.607 1.588 1.755 1.755 1.848 1.927 1.766 1.750	6.92
79) S	Terphenyl-d14	1.101 1.089 1.200 1.214 1.287 1.328 1.232 1.207	7.30
80)	Butylbenzylphth...	0.523 0.568 0.692 0.671 0.685 0.716 0.682 0.648	11.20
81)	Benzo(a)anthra...	1.292 1.270 1.374 1.309 1.338 1.394 1.336 1.330	3.31
82)	3,3'-Dichlorob...	0.319 0.336 0.386 0.367 0.371 0.394 0.362 0.362	7.31
83)	Chrysene	1.237 1.229 1.321 1.234 1.249 1.321 1.252 1.263	3.18
84)	Bis(2-ethylhex...	0.583 0.632 0.799 0.757 0.777 0.827 0.822 0.742	12.95
85) c	Di-n-octyl pht...	0.801 1.049 1.119 1.188 1.342 1.315 1.136	17.49

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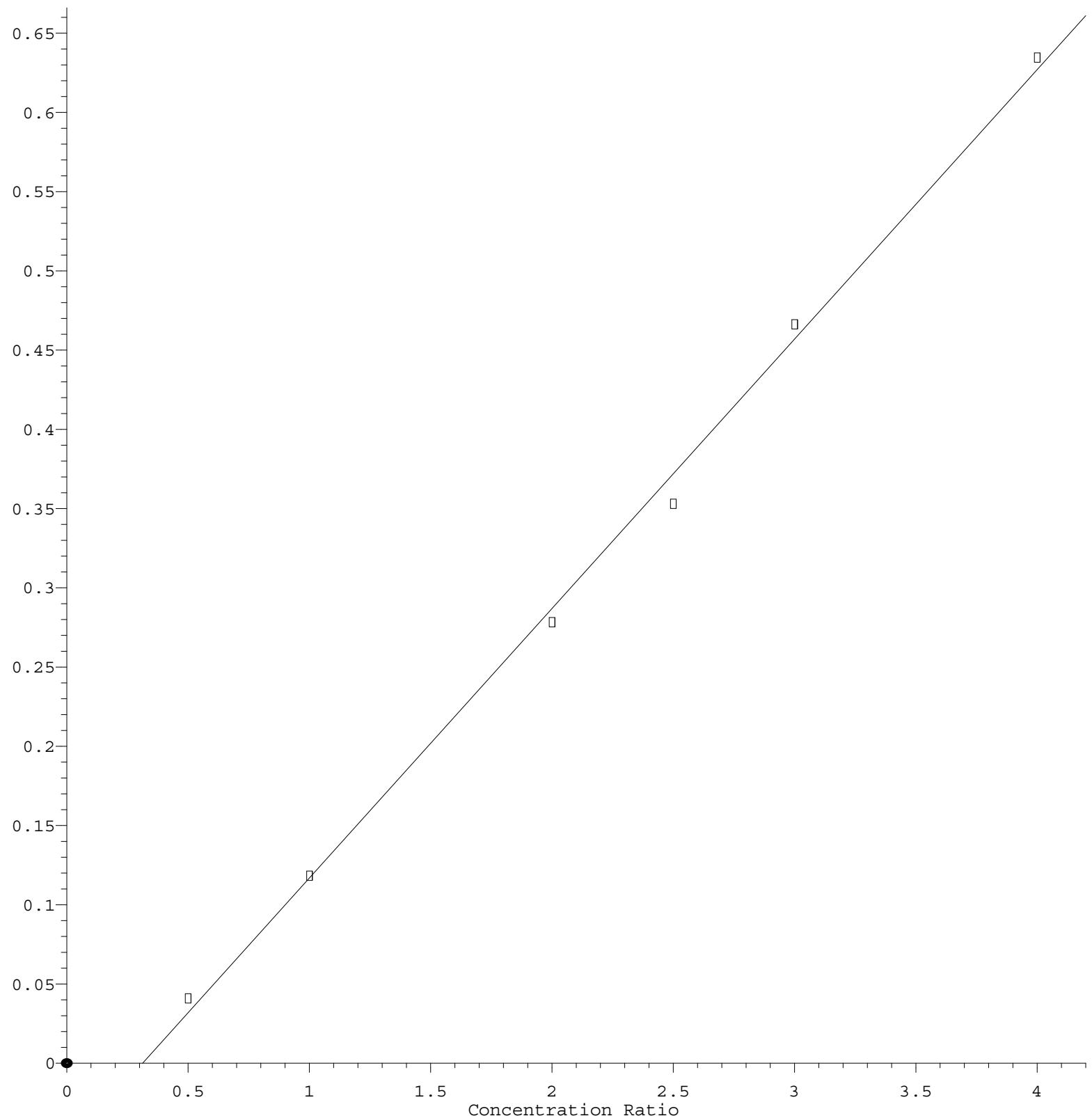
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86)	I	Perylene-d12	-----ISTD-----								
87)		Indeno(1,2,3-c...)	1.098	1.179	1.322	1.474	1.578	1.680	1.598	1.418	15.72
88)		Benzo(b)fluora...	1.287	1.291	1.357	1.301	1.286	1.295	1.321	1.305	1.97
89)		Benzo(k)fluora...	1.286	1.265	1.375	1.273	1.234	1.315	1.241	1.284	3.77
90)	C	Benzo(a)pyrene	0.979	0.947	1.043	1.043	1.047	1.106	1.074	1.034	5.23
91)		Dibenzo(a,h)an...	0.896	0.968	1.085	1.200	1.293	1.377	1.326	1.163	15.94
92)		Benzo(g,h,i)pe...	0.914	0.976	1.084	1.214	1.313	1.379	1.323	1.172	15.58

(#) = Out of Range

2,4-Dinitrophenol

Response Ratio



$$\text{Response} = 1.700\text{e-001} * \text{Amt} - 5.317\text{e-002}$$

Coef of Det (r^2) = 0.997269 Curve Fit: Linear

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Calibration Table Last Updated: Wed Sep 14 14:55:51 2022